

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094401.D
 Acq On : 15 Oct 2017 19:59
 Operator : SJ/JU
 Sample : I5548-22DL 5X
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE8DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:26:35 PM

Quant Time: Oct 16 08:07:48 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 07:45:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1975	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	10301	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5947	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12668	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12904	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	11948	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	731	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1706	0.04	ng/ul	0.00
Target Compounds						
					Ovalue	
3) Naphthalene	10.75	128	1536	0.061	ng/ul#	89
5) 2-Methylnaphthalene	12.36	142	1499	0.092	ng/ul	97
7) Acenaphthylene	14.25	152	2263	0.098	ng/ul#	92
8) Acenaphthene	14.59	153	1509	0.079	ng/ul	99
9) Fluorene	15.57	166	1941	0.087	ng/ul	90
12) Phenanthrene	17.31	178	41674	1.259	ng/ul	95
13) Anthracene	17.39	178	11206	0.348	ng/ul	96
15) Fluoranthene	19.32	202	110424	2.539	ng/ul	84
17) Pyrene	19.68	202	101565	2.829	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	62793	1.834	ng/ul#	88
19) Chrysene	21.47	228	58879m	1.489	ng/ul	
21) Benzo(b)fluoranthene	23.19	252	87180m	2.686	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	21752m	0.580	ng/ul	
23) Benzo(a)pyrene	23.85	252	57763	1.713	ng/ul	90
24) Indeno(1,2,3-cd)pyrene	26.65	276	40062	1.161	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.65	278	10153	0.351	ng/ul#	40
26) Benzo(a,h,i)perylene	27.50	276	40359	1.356	ng/ul#	88

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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